#### Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

#### Listing of Claims:

1. (Currently amended) A compound for modulating kinase activity of Formula I,



or a pharmaceutically acceptable salt thereof, wherein,

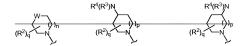
Ar is selected from the following formulae

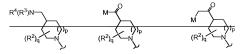


wherein Ar is substituted with —X and -Y-L-Z, in an ortho relationship to each other, and said five—to-six-membered aromatic ring system Ar is optionally substituted with up to four R<sup>1</sup>:

each  $R^1$  is independently selected from -H, halogen, -CN, -NO<sub>2</sub>, -OR<sup>3</sup>, -N( $R^3$ )R<sup>3</sup>, -S(O)<sub>0-2</sub>R<sup>3</sup>, -SO<sub>2</sub>N(R<sup>3</sup>)R<sup>3</sup>, -CO<sub>2</sub>R<sup>3</sup>, -C(O)N(R<sup>3</sup>)R<sup>3</sup>, -N(R<sup>3</sup>)SO<sub>2</sub>R<sup>3</sup>, -N(R<sup>3</sup>)C(O)R<sup>3</sup>, -N(R<sup>3</sup>)CO<sub>2</sub>R<sup>3</sup>, -C(O)R<sup>3</sup>, -OC(O)R<sup>3</sup>, optionally substituted lower alkyl, optionally substituted aryl, optionally substituted lower arylalkyl, optionally substituted heterocyclyl, and optionally substituted lower heterocyclylalkyl:

X is selected from the following six formulae:





wherein.

W is selected from  $-C(R^2)(R^2)$ ,  $-N(R^4)$ ,  $-S(O)_{0-2}$ , and -O;

X is selected from the following formulae.

# wherein R<sup>4a</sup> is -C(O)N(R<sup>3</sup>)R<sup>3</sup>:

n = 1 or 2;

p = 0 or 1;

q is 1 to 3;

#### M is OR3 or N(R3)R4:

each  $R^2$  is independently selected from -H, halogen, oxo, -CN, -NH<sub>2</sub>, -NO<sub>2</sub>, -OR<sup>3</sup>, -N(R<sup>3</sup>)R<sup>3</sup>, -N(R<sup>3</sup>)R<sup>5</sup>, -S(O) $_0$ 2R<sup>3</sup>, -SO<sub>2</sub>N(R<sup>3</sup>)R<sup>3</sup>, -CO<sub>2</sub>R<sup>3</sup>, -C(O)N(R<sup>3</sup>)R<sup>3</sup>, -N(R<sup>3</sup>)SO<sub>2</sub>R<sup>3</sup>, -N(R<sup>3</sup>)C(O)R<sup>3</sup>, -N(R<sup>3</sup>)CO<sub>2</sub>R<sup>3</sup>, -N(R<sup>3</sup>)C(O)N(R<sup>3</sup>)R<sup>3</sup>, -C(O)R<sup>3</sup>, -OC(O)R<sup>3</sup>, optionally substituted lower alkyl, optionally substituted aryl, optionally substituted lower arylalkyl, optionally substituted heterocyclyl, and optionally substituted lower heterocyclylalkyl;

two of R<sup>2</sup>, together with the atoms to which they are attached, can form an optionally substituted three- to seven-membered ring system;

each R<sup>3</sup> is independently selected from -H, optionally substituted lower alkyl, optionally substituted aryl, optionally substituted lower arylalkyl, optionally substituted heterocyclyl, and optionally substituted lower heterocyclylalkyl; or

two of R<sup>3</sup>, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl ring, said optionally substituted five- to seven-membered heterocyclyl ring optionally containing at least one additional heteroatom selected from N, O, S, and P;

each  $R^4$  is independently selected from  $R^3$ ,  $-SO_2R^3$ ,  $-SO_2N(R^3)R^3$ ,  $-CO_2R^3$ ,  $-C(O)N(R^3)R^3$ , and  $-C(O)R^3$ ;

-Y-L-Z is selected from the following formulae,

wherein g is zero to two; D is selected from  $-C(R^5)(R^5)$ -, -O-,  $-S(O)_{0\cdot 2^-}$ , and  $-N(R^4)$ -; Q is =N- or  $-C(R^5)$ -, T is selected from absent,  $-N(R^3)$ -, -S- and -O-; and each methylene between Y and T is optionally substituted; provided that when both Y and T are heteroatoms then g must be two;

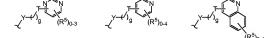
Y is selected from -CH2-, -O-, -S(O)0-2-, -N(R3)-, and absent;

 $R^5$  is selected from -H, halogen, -CN, -NO<sub>2</sub>, -OR<sup>3</sup>, -N(R<sup>3</sup>)R<sup>4</sup>, -S(O)<sub>0.2</sub>R<sup>3</sup>, -SO<sub>2</sub>N(R<sup>3</sup>)R<sup>3</sup>, -CO<sub>2</sub>R<sup>3</sup>, -C(O)N(R<sup>3</sup>)R, -N(R<sup>3</sup>)SO<sub>2</sub>R<sup>3</sup>, -N(R<sup>3</sup>)C(O)R<sup>3</sup>, -N(R<sup>3</sup>)CO<sub>2</sub>R<sup>3</sup>, -C(O)R<sup>3</sup>, optionally substituted lower alkyl, optionally substituted aryl, optionally substituted lower arylalkyl, optionally substituted heterocyclyl, and optionally substituted lower heterocyclylalkyl; and

optionally two of  $R^5$ , together with the atoms to which they are attached, form a second ring system fused with said five- to seven-membered ring system, said second ring system substituted with zero to four of  $R^5$ .

- (Canceled)
- (Canceled)
- (Canceled)

- (currently amended) The compound according to claim 4 <u>1</u>, wherein Y is -O- or optionally substituted -CH<sub>2</sub>-.
- (Canceled)
- (Canceled)
- (original) The compound according to claim 7 <u>1</u>, wherein each R<sup>2</sup> is independently selected from -H or optionally substituted lower alkyl.
- (original) The compound according to claim 8, wherein each R<sup>2</sup> is independently selected from -H, haloalkyl, -C<sub>1-6</sub>alkyl-N(R<sup>3</sup>)R<sup>3</sup>, -C<sub>1-6</sub>alkyl-OR<sup>3</sup>, -C<sub>1-6</sub>alkyl-CO<sub>2</sub>R<sup>3</sup>, and -C<sub>1-6</sub>alkyl-C(O)N(R<sup>3</sup>)R<sup>3</sup>.8
- 10. (Canceled)
- (currently amended) The compound according to claim 9 +0, wherein -Y-L-Z is selected from the following formulae,



wherein Y, T, and g are as described above.

- 12. (original) The compound according to claim 11, wherein g is one or two.
- 13. (original) The compound according to claim 12, wherein each R<sup>5</sup> is independently selected from -H, halogen, -CN, -NH<sub>2</sub>, -NO<sub>2</sub>, -OR<sup>3</sup>, -N(R<sup>3</sup>)R<sup>4</sup>, -S(O)<sub>0-2</sub>R<sup>3</sup>, -SO<sub>2</sub>N(R<sup>3</sup>)R<sup>3</sup>, -CO<sub>2</sub>R<sup>3</sup>, -C(O)N(R<sup>3</sup>)R<sup>3</sup>, -N(R<sup>3</sup>)SO<sub>2</sub>R<sup>3</sup>, -N(R<sup>3</sup>)C(O)R<sup>3</sup>, -N(R<sup>3</sup>)CO<sub>2</sub>R<sup>3</sup>, -C(O)R<sup>3</sup>, and optionally substituted lower alkyl.
- (original) The compound according to claim 13, wherein -Y-L-Z is selected from the following formulae.

$$R^3$$
  $R^{4b}$   $R^3$   $R^4$ 

15. (currently amended) The compound according to claim 14, having formula III,

ш

wherein J is N or CH, and B is =N- or = $C(R^5)$ -.

- (original) The compound according to claim 15, wherein R<sup>3a</sup> is selected from optionally substituted aryl, optionally substituted lower arylalkyl, optionally substituted heterocyclyl, and optionally substituted lower heterocyclylalkyl.
- (original) The compound according to claim 16, wherein R<sup>3a</sup> is selected from optionally substituted aryl and optionally substituted heteroaryl.
- (original) The compound according to claim 17, wherein R<sup>3a</sup> is optionally substituted phenyl.
- 19. (original) The compound according to claim 18, wherein said optionally substituted phenyl is substituted with at least one of halogen, -CN, -CF<sub>3</sub>, -NH<sub>2</sub>, -NO<sub>2</sub>, -OR<sup>3</sup>, -N(R<sup>3</sup>)R<sup>3</sup>, -S(O)<sub>0-2</sub>R<sup>3</sup>, -SO<sub>2</sub>N(R<sup>3</sup>)R<sup>3</sup>, -CO<sub>2</sub>R<sup>3</sup>, -CO<sub>2</sub>R<sup>3</sup>, -CO<sub>3</sub>N<sup>3</sup>, -N(R<sup>3</sup>)SO<sub>2</sub>R<sup>3</sup>, -N(R<sup>3</sup>)CO<sub>2</sub>R<sup>3</sup>, -N(R<sup>3</sup>)CO<sub>2</sub>R<sup>3</sup>, -C(O)R<sup>3</sup>, optionally substituted lower alkyl, and optionally substituted aryl.
- (original) The compound according to claim 19, wherein said optionally substituted
  phenyl group is substituted with at least one trifluoromethyl group.

- (original) The compound according to claim 20, wherein said optionally substituted
  phenyl group is substituted with at least two trifluoromethyl groups
- (original) The compound according to claim 19, wherein said optionally substituted
  phenyl group is substituted with at least one lower alkyl group.
- 23. (original) The compound according to claim 19, wherein R3b is -H.
- (original) The compound according to claim 23, wherein R<sup>4b</sup> is selected from R<sup>3</sup>, -H, -CO<sub>2</sub>R<sup>3</sup>, -C(O)N(R<sup>3</sup>)R<sup>4</sup>, and -C(O)R<sup>3</sup>.
- (Canceled)
- (Canceled)
- (original) The compound according to claim 24, wherein Ar is according to the formula below.

 (original) The compound according to claim 24, wherein Ar is according to the formula below.

$$\gtrsim$$

- (Canceled)
- 30. (Currently amended) A compound for modulating kinase activity of Formula IV,

$$(R^{6})_{0.5} \xrightarrow{(R^{2})_{0.4}} \xrightarrow{B} \overset{H}{\overset{H}{\overset{}}}$$

or a pharmaceutically acceptable salt thereof, wherein,

Ar is selected from the following formulae:

$$(R^{1})_{0\cdot 2} \qquad (R^{1})_{0\cdot 3} \qquad (R^{1})_{0\cdot 3} \qquad (R^{1})_{0\cdot 3} \qquad (R^{1})_{0\cdot 4}$$

each  $R^1$  is independently selected from -H, halogen, -CN, -NO<sub>2</sub>, -OR<sup>3</sup>, -N( $R^3$ )R<sup>3</sup>, -S(O)<sub>6-2</sub>R<sup>3</sup>, -SO<sub>2</sub>N(R<sup>3</sup>)R<sup>3</sup>, -CO<sub>2</sub>R<sup>3</sup>, -C(O)N(R<sup>3</sup>)R<sup>3</sup>, -N(R<sup>3</sup>)SO<sub>2</sub>R<sup>3</sup>, -N(R<sup>3</sup>)CO)R<sup>3</sup>, -N(R<sup>3</sup>)CO<sub>2</sub>R<sup>3</sup>, -C(O)R<sup>3</sup>, optionally substituted lower alkyl, optionally substituted aryl, optionally substituted lower arylalkyl, optionally substituted heterocyclyl, and optionally substituted lower heterocyclylalkyl;

optionally two of  $\mathbb{R}^1$ , together with the atoms to which they are attached, form a first ring system fused with Ar, said first ring system substituted with zero to three additional of  $\mathbb{R}^1$ .

each  $R^2$  is independently selected from -H, halogen, oxo, -CN, -NH<sub>2</sub>, -NO<sub>2</sub>, -OR<sup>3</sup>, -N(R<sup>3</sup>)R<sup>3</sup>, -N(R<sup>3</sup>)R<sup>5</sup>, -S(O) $_{0}$ -R<sup>3</sup>, -SO<sub>2</sub>N(R<sup>3</sup>)R<sup>3</sup>, -CO<sub>2</sub>R<sup>3</sup>, -C(O)N(R<sup>3</sup>)R<sup>3</sup>, -N(R<sup>3</sup>)SO<sub>2</sub>R<sup>3</sup>, -N(R<sup>3</sup>)C(O)R<sup>3</sup>, -N(R<sup>3</sup>)C(O)R<sup>3</sup>, -N(R<sup>3</sup>)C(O)R<sup>3</sup>, -C(O)R<sup>3</sup>, optionally substituted lower alkyl, optionally substituted aryl, optionally substituted lower arylalkyl, optionally substituted heterocyclyl, and optionally substituted lower heterocyclylalkyl;

two of R<sup>2</sup>, together with the atoms to which they are attached, can form an optionally substituted three- to seven-membered ring system;

each R<sup>3</sup> is independently selected from -H, optionally substituted lower alkyl, optionally substituted aryl, optionally substituted lower arylalkyl, optionally substituted heterocyclyl, and optionally substituted lower heterocyclylalkyl; or

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two of R<sup>3</sup>, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl ring, said optionally substituted five- to seven-membered heterocyclyl ring optionally containing at least one additional heteroatom selected from N, O, S, and P;

each  $R^4$  is independently selected from  $R^3$ ,  $-SO_2R^3$ ,  $-SO_2N(R^3)R^3$ ,  $-CO_2R^3$ ,  $-C(O)N(R^3)R^3$ , and  $-C(O)R^3$ ;

Y is selected from optionally substituted -CH<sub>2</sub>-, -O-, -S-, and -N(R<sup>3</sup>)-;

L is selected from optionally substituted -CH2-, -O-, -S-, -N(R $^3$ )- and absent;

provided that Y and L are not both heteroatoms;

B is =N- or =C(H)-;

at each instance,  $R^3$  and  $R^6$  are independently selected from -H, halogen, -CN, -NO<sub>2</sub>, -OR<sup>3</sup>, -N(R<sup>3</sup>)R<sup>4</sup>, -S(O)<sub>0-2</sub>R<sup>3</sup>, -SO<sub>2</sub>N(R<sup>3</sup>)R<sup>3</sup>, -CO<sub>2</sub>R<sup>3</sup>, -CO<sub>2</sub>N(R<sup>3</sup>)R, -N(R<sup>3</sup>)SO<sub>2</sub>R<sup>3</sup>, -N(R<sup>3</sup>)C(O)R<sup>3</sup>, -N(R<sup>3</sup>)CO<sub>2</sub>R<sup>3</sup>, -C(O)R<sup>3</sup>, optionally substituted aryl, optionally substituted aryl, optionally substituted lower arylalkyl, optionally substituted heterocyclyl, and optionally substituted lower heterocyclylalkyl; and

optionally two of R<sup>5</sup>, together with the atoms to which they are attached, form a ring system fused with the ring containing B according to formula IV, said ring system substituted with zero to two additional of R<sup>5</sup>.

- (original) The compound according to claim 30, wherein Y is -O- and L is optionally substituted -CH<sub>2</sub>-.
- (original) The compound according to claim 31, wherein at least one of R<sup>6</sup> is optionally substituted lower alkyl.
- (original) The compound according to claim 32, wherein said at least one optionally substituted lower alkyl is meta- to the piperazine urea function as depicted in formula IV.
- (original) The compound according to claim 33, wherein R<sup>4a</sup> is selected from R<sup>3</sup>, -H, -CO<sub>2</sub>R<sup>3</sup>, -C(O<sub>2</sub>N(R<sup>3</sup>)R<sup>4</sup>, and -C(O<sub>2</sub>R<sup>3</sup>.

- (original) The compound according to claim 34, wherein R<sup>4a</sup> is selected from -H, -CO<sub>2</sub>R<sup>3</sup>,
  -C(O)N(R<sup>3</sup>)R<sup>4</sup>, and -C(O)R<sup>3</sup>.
- 36. (original) The compound according to claim 35, wherein -Y-L- is -OCH<sub>2</sub>-.
- 37. (Canceled)
- (Canceled)
- (original) The compound according to claim 36, wherein Ar is according to the formula below.

$$\mathbb{Z}$$

 (original) The compound according to claim 36, wherein Ar is according to the formula below.

$$\sum_{i}$$

- 41. (Cancelled)
- 42. The compound according to claim 1, A compounds selected from Table 4.

Table 4

97	N-[3,5-bis(trifluoromethyl)phenyl]-4- {3-[(pyridin-4-ylmethyl)oxy]pyridin-2- yl}piperazine-1-carboxamide	HN N N N N N N N N N N N N N N N N N N
103	N-(4-chlorophenyl)-4-{3-[(pyridin-4-ylmethyl)oxy]pyridin-2-yl}piperazine- 1-carboxamide	

105	N-(3-chlorophenyl)-4-{3-[(pyridin-4- ylmethyl)oxy]pyridin-2-yl}piperazine- 1-carboxamide	
142	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-[3,5- bis(trifluoromethyl)phenyl]piperazine-1- carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N
144	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-(3- ethylphenyl)piperazine-1-carboxamide	N N N N N N N N N N N N N N N N N N N
161	methyl [4-({[2-(4-{[(3- ethylphenyl)amino carbonyl}piperazin- 1-yl)pyridin-3-yl]oxy}methyl)pyridin-2- yl]carbamate	
164	methyl [4-({[2-(4-{[(3- bromophenyl)amino]carbonyl}piperazin -1-yl)pyridin-3-yl]oxy}methyl)pyridin- 2-yl]carbamate	NH NH NH
165	methyl {4-[({2-[4-({[3- (methyloxy)phenyl]amino;carbonyl)pip erazin-1-yl]pyridin-3- yl}oxy)methyl]pyridin-2-yl}carbamate	

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Table 4

166	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy] pyridin-2-yl)-N-[3- (methyloxy)phenyl]piperazine-1- carboxamide	N N N N N N N N N N N N N N N N N N N
167	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy} pyridin-2-yl)-N-[3-(1- methylethyl)phenyl]piperazine-1- carboxamide	NH <sub>2</sub> N N N N N N N N N N N N N N N N N N N
168	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-{3- [(trifluoromethyl)oxy]phenyl}piperazine -1-carboxamide	NH2 N N N N N N N N N N N N N N N N N N
169	4-(3-[[(2-aminopyrimidin-4- yl)methyl]oxy]pyridin-2-yl)-N-[2- fluoro-5- (trifluoromethyl)phenyl]piperazine-1- carboxamide	NH <sub>2</sub>
170	N-(3-ethylphenyl)-4-[3-({[2-({[(3-ethylphenyl)amino]earbonyl]amino)pyridin-4-yl]methyl}oxy)pyridin-2-yl]piperazine-1-earboxamide	HN NH NH

Table 4

171	N-(3-ethylphenyl)-4-(3-{[(2-{[(4- methylpiperazin-1- yl)acetyl]amino}pyridin-4- yl)methyl]oxy}pyridin-2-yl)piperazine- I-carboxamide	N N N N N N N N N N N N N N N N N N N
173	N-[3,5-bis(trifluoromethyl)phenyl]-4-(3- {[(2-{[(4-methylpiperazin-1- yl)acetyl]amino}pyridin-4- yl)methyl]oxy}pyridin-2-yl)piperazine- l-carboxamide	N N N N N N N N N N N N N N N N N N N
174	4-(3-{{(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-[3- (trifluoromethyl)phenyl]piperazine-1- carboxamide	H <sub>2</sub> N N N NH
175	4-[3-({[2-(acetylamino)pyridin-4-yl]methyl}oxy)pyridin-2-yl]-N-(3-ethylphenyl)piperazine-1-carboxamide	NH NH NH NH
176	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-(3-ethyl- 4-fluorophenyl)piperazine-1- carboxamide	H <sub>2</sub> N N N NH

Table 4

177	2-[4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)piperazin-1- yl]-N-[3,5- bis(trifluoromethyl)phenyl]acetamide	N N N N N N N N N N N N N N N N N N N
178	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N- phenylpiperazine-1-carboxamide	H <sub>2</sub> N NH
179	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy  pyridin-2-yl)-N-(3- chloro-5-ethylphenyl)piperazine-1- carboxamide	H <sub>2</sub> N N N CI
180	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-(5-ethyl- 2-fluorophenyl)piperazine-1- carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N
181	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy] pyridin-2-yl)-N-(3- bromo-5-ethylphenyl)piperazine-1- carboxamide	H <sub>2</sub> N N N NH

Table 4

182	2-(4-methylpiperazin-1-yl)ethyl [4-({[2- (4-{[(3- ethylphenyl)amino]carbonyl}piperazin- 1-yl)pyridin-3-yl]oxyl methyl)pyridin-2- yl]carbamate	
183	4-(3-{[(2-aminopyrimidin-4-yl)methyl]oxy}pyridin-2-yl)-N-(3-chlorophenyl)piperazine-1-carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N
184	4-(3-{[(2-aminopyrimidin-4-yl)methyl]oxy}pyridin-2-yl)-N-(3-bromophenyl)piperazine-1-carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N
185	N-[4-({[2-(4-acetylpiperazin-1-yl)pyridin-3-yl]oxy}methyl)pyridin-2-yl]-2-(4-methylpiperazin-1-yl)acetamide	
186	4-(3-{[(2-aminopyrimidin-4-yl)methyl]oxy}pyridin-2-yl)-N-(3-fluorophenyl)piperazine-1-carboxamide	H <sub>2</sub> N N O NH
187	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-(4- fluorophenyl)piperazine-1-carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N

Table 4

	,	
188	4-(3-{[(2-aminopyrimidin-4-yl)methyl]oxy}pyridin-2-yl)-N-(2-fluorophenyl)piperazine-1-carboxamide	H <sub>2</sub> N NH NH
189	4-(3-{[(2-aminopyrimidin-4-yl)methyl]oxy}pyridin-2-yl)-N-(3,5-diethylphenyl)piperazine-1-carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N
190	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}-5-bromopyridin-2-yl)- N-(3-ethylphenyl)piperazine-1- carboxamide	Br N N NH
191	N-methyl-4-(3-{[(2-{[(4- methylpiperazin-1- yl)acetyl]amino}pyridin-4- yl)methyl]oxy}pyridin-2-yl)piperazine- 1-carboxamide	NH NH NH
192	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-[2- chloro-5- (trifluoromethyl)phenyl]piperazine-1- carboxamide	H <sub>2</sub> N N F <sub>3</sub> C CI
193	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy pyridin-2-yl)-N-(5- chloro-2-fluorophenyl)piperazine-1- carboxamide	H <sub>2</sub> N N CI F
194	4-(3-{[(2-amino-5-bromopyrimidin-4-yl)methyl]oxy]-5-bromopyridin-2-yl)-N-(3-ethylphenyl)piperazine-1-carboxamide	Br N N N N N N N N N N N N N N N N N N N

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Table 4

195	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-[2- fluoro-3- (trifluoromethyl)phenyl]piperazine-1- carboxamide	H <sub>2</sub> N NH NH CF <sub>3</sub>
196	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-[3- fluoro-5- (trifluoromethyl)phenyl]piperazine-1- carboxamide	H <sub>2</sub> N NH F-CF <sub>3</sub>
197	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-(3,5- dichlorophenyl)piperazine-1- carboxamide	H <sub>2</sub> N N CI CI
198	N-(3-chloro-5-ethylphenyl)-4-(3-{[(2- {[(4-methylpiperazin-1- yl)acetyl]amino}pyridin-4- yl)methyl]oxy}pyridin-2-yl)piperazine- 1-carboxamide	N N N N N N N N N N N N N N N N N N N
199	N-(5-ethyl-2-fluorophenyl)-4-(3-{[(2- {[(4-methylpiperazin-1- yl)acetyl]amino}pyridin-4- yl)methyl]oxy}pyridin-2-yl)piperazine- 1-carboxamide	NH NH NH

Table 4

200	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-[3-ethyl- 5-(trifluoromethyl)phenyl]piperazine-1- carboxamide	H <sub>2</sub> N NH NH CF <sub>3</sub>
204	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N- methylpiperazine-1-carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N
205	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N- ethylpiperazine-1-carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N
206	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N- cyclohexylpiperazine-1-carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N
207	4-({[2-(4-acctylpiperazin-1-yl)pyridin-3-yl]oxy}methyl)pyrimidin-2-amine	
208	4-{{[2-(4-propanoylpiperazin-1-yl)pyridin-3-yl]oxy}methyl)pyrimidin- 2-amine	○ N N N N N N N N N N N N N N N N N N N
209	N-(3-cyclopropylphenyl)-4-(3-{[(2-{[(4-methylpiperazin-1-yl)acetyl]amino}pyridin-4-yl)methyl]oxy}pyridin-2-yl)piperazine-l-carboxamide	NH NH

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210	4-(3-[[(2-aminopyrimidin-4- yl)methyl]oxy] pyridin-2-yl)-N-(3- cyclopropylphenyl)piperazine-1- carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N
211	N-[2-fluoro-5-(trifluoromethyl)phenyl]- 4-(3-{[(2-{[(4-methylpiperazin-1- yl)acetyl]amino}pyridin-4- yl)methyl]oxy}pyridin-2-yl)piperazine- l-carboxamide	N N N N N N N N N N N N N N N N N N N
212	N-[3-fluoro-5-(trifluoromethyl)phenyl]- 4-(3-{[(2-{[(4-methylpiperazin-1- yl)acetyl]amino}pyridin-4- yl)methyl]oxy}pyridin-2-yl)piperazine- l-carboxamide	N N N N N N N N N N N N N N N N N N N
213	N-(3,5-dichlorophenyl)-4-(3-{[(2-{[(4-methylpiperazin-1-yl)acetyl]amino}pyridin-4-yl)methyl]oxy}pyridin-2-yl)piperazine-l-carboxamide	N N N N N N N N N N N N N N N N N N N
214	4-(3-{[(2-{[(4-methylpiperazin-1-yl)acetyl]amino}pyridin-4-yl)methyl]oxy}pyridin-2-yl)-N-[3-(trifluoromethyl)phenyl]piperazine-1-carboxamide	N— N— N— N— NH NH NH NH

216	4-(3-{[1-(2-aminopyrimidin-4- yl)ethyl]oxy] pyridin-2-yl)-N-[3,5- bis(trifluoromethyl)phenyl]piperazine-1- carboxamide	$H_2N$ $N$ $F_3C$ $CF_3$
219	4-[{2-[4-(3,4-dihydroquinolin-1(2H)- ylcarbonyl)piperazin-1-yl]pyridin-3- yl}oxy)methyl]pyrimidin-2-amine	NH <sub>2</sub>
220	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-(2- methylpropyl)piperazine-1-carboxamide	H <sub>2</sub> N N N NH
226	N-(3,5-diethylphenyl)-4-(3-{[(2-{[(4- methylpiperazin-1- yl)acetyl]amino]pyridin-4- yl)oxy]methyl]pyridin-2-yl)piperazine- 1-carboxamide	NH NH
227	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}-6-methylpyridin-2-yl)- N-(3-ethylphenyl)piperazine-1- carboxamide	H <sub>2</sub> N NH
228	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}-6-methylpyridin-2-yl)- N-[3,5- bis(trifluoromethyl)phenyl]piperazine-1- carboxamide	$H_2N$ $F_3C$ $CF_3$

233	N-[3-chloro-5-(trifluoromethyl)phenyl]- 4-(3-{[(2-{[(4-methylpiperazin-1- yl)acetyl]amino}pyridin-4- yl)methyl]oxy}pyridin-2-yl)piperazine- l-carboxamide	N N N N N N N N N N N N N N N N N N N
235	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-[3- chloro-5- (trifluoromethyl)phenyl]piperazine-1- carboxamide	H <sub>2</sub> N N H NH CI
237	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}-6-chloropyridin-2-yl)- N-(3-ethylphenyl)piperazine-1- carboxamide	CI N N N N N N N N N N N N N N N N N N N
243	N-[3-chloro-2-fluoro-5- (trifluoromethyl)phenyl]-4-(3-{[(2-{[(4- methylpiperazin-1- yl)acetyl]amino}pyridin-4- yl)methyl]oxy}pyridin-2-yl)piperazine- 1-carboxamide	NH NH F <sub>3</sub> C — F
244	4-(3-{[(2-aminopyrimidin-4-yl)methyl]oxy}-6-chloropyridin-2-yl)-N-[3,5-bis(trifluoromethyl)phenyl]piperazine-1-carboxamide	$\begin{array}{c c} CI & & \\ $

Table 4

245	4-(3-{[1-(2-aminopyrimidin-4-yl)ethyl]oxy}pyridin-2-yl)-N-(3-ethylphenyl)piperazine-1-carboxamide	H <sub>2</sub> N N NH
246	4-(3-{[(2-aminopyrimidin-4-yl)methyl]oxy}-6-chloropyridin-2-yl)-N-(5-ethyl-2-fluorophenyl)piperazine-1-carboxamide	H <sub>2</sub> N N N NH
247	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-(3-ethyl- 5-fluorophenyl)piperazine-1- carboxamide	H <sub>2</sub> N N NH
249	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-[3- chloro-2-fluoro-5- (trifluoromethyl)phenyl]piperazine-1- carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N
250	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-[3,5- bis(trifluoromethyl)phenyl]-N- methylpiperazine-1-carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N

- 43. A pharmaceutical composition comprising the compound according to claim 1 and a pharmaceutically acceptable carrier.
- 44. (Canceled).
- 45. (Canceled)
- 46. (Canceled)
- 47. (Canceled)

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- 48. (Canceled)
- 49. (Canceled)
- 50. (Canceled).
- 51. (Canceled)
- 52. (Canceled)
- 53. (Canceled)
- 54. (Canceled)
- 55. (Canceled)
- 56. (Canceled)
- 57. (Canceled)